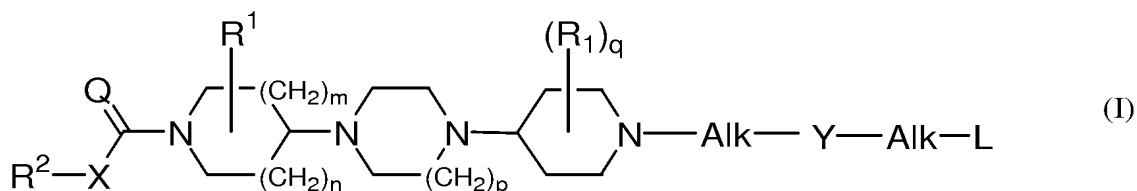


This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims:

1. (Previously Presented/Withdrawn) A compound according to the general Formula (I)



the pharmaceutically acceptable acid or base addition salts thereof, the stereochemically isomeric forms thereof, the *N*-oxide form thereof and prodrugs thereof, wherein :

- n is an integer, equal to 0, 1 or 2 ;
m is an integer, equal to 1 or 2, provided that if m is 2, then n is 1 ;
p is an integer equal to 1 or 2 ;
Q is O or NR³ ;
X is a covalent bond or a bivalent radical of formula -O-, -S- or -NR³- ;
each R³ independently from each other, is hydrogen or alkyl ;
each R¹ independently from each other, is selected from the group of Ar¹, Ar¹-alkyl and di(Ar¹)-alkyl ;
q is an integer equal to 0 or 1 ;
R² is selected from the group consisting of alkyl, Ar², Ar²-alkyl, Het¹ or Het¹-alkyl ;
Y is a covalent bond or a bivalent radical of formula -C(=O)- or -SO₂-;
each Alk represents, independently from each other, a covalent bond; a bivalent straight or branched, saturated or unsaturated hydrocarbon radical having from 1 to 6 carbon atoms ; or a cyclic saturated or unsaturated hydrocarbon radical having from 3 to 6 carbon atoms ; each radical optionally substituted on one or more carbon atoms with one or more alkyl, phenyl, halo, cyano, hydroxy, formyl and amino radicals ;
L is selected from the group of hydrogen, alkyloxy, Ar³-oxy, alkyloxycarbonyl, mono- and di(alkyl)amino, mono- and di(Ar³)amino, mono- and di(alkyloxycarbonyl)amino, Ar³, Ar³-carbonyl, Het² and Het²-carbonyl;

- Ar¹ is phenyl, optionally substituted with 1, 2 or 3 substituents each independently from each other selected from the group consisting of halo, alkyl, cyano, aminocarbonyl and alkyloxy ;
- Ar² is naphthalenyl or phenyl, each optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group consisting of halo, nitro, amino, mono- and di(alkyl)amino, cyano, alkyl, hydroxy, alkyloxy, carboxyl, alkyloxycarbonyl, aminocarbonyl and mono- and di(alkyl)aminocarbonyl ;
- Ar³ is naphthalenyl or phenyl, optionally substituted with 1, 2 or 3 substituents each independently from each other selected from the group consisting of alkyloxy, alkyl, halo, hydroxy, pyridinyl, morpholinyl, pyrrolidinyl, imidazo[1,2-*a*]pyridinyl, morpholinylcarbonyl, pyrrolidinylcarbonyl, amino and cyano;
- Het¹ is a monocyclic heterocyclic radical selected from the group consisting of pyrrolyl, pyrazolyl, imidazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, thiadiazolyl, isothiazolyl, pyridinyl, pyrimidinyl, pyrazinyl and pyridazinyl ; or a bicyclic heterocyclic radical selected from the group of quinolinyl, quinoxalinyl, indolyl, benzimidazolyl, benzoxazolyl, benzisoxazolyl, benzothiazolyl, benzisothiazolyl, benzofuranyl and benzothienyl ; each heterocyclic radical may optionally be substituted on any atom by a radical selected from the group consisting of halo and alkyl ;
- Het² is a monocyclic heterocyclic radical selected from the group consisting of pyrrolidinyl, dioxolyl, imidazolidinyl, pyrazolidinyl, piperidinyl, morpholinyl, dithianyl, thiomorpholinyl, piperazinyl, imidazolidinyl, tetrahydrofuranyl, 2H-pyrrolyl, pyrrolinyl, imidazolinyl, pyrrazolinyl, pyrrolyl, imidazolyl, pyrazolyl, triazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, thiadiazolyl, isothiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyridazinyl and triazinyl ; or a bicyclic heterocyclic radical selected from the group consisting of benzopiperidinyl, quinolinyl, quinoxalinyl, indolyl, isoindolyl, chromenyl, benzimidazolyl, imidazo[1,2-*a*]pyridinyl, benzoxazolyl, benzisoxazolyl, benzothiazolyl, benzisothiazolyl, benzofuranyl and benzothienyl ; each radical optionally substituted with one or more radicals selected from the group consisting of Ar¹, Ar¹alkyl, halo, hydroxy, alkyl, piperidinyl, pyrrolyl, thienyl, oxo, alkyloxy, alkyloxyalkyl and alkyloxycarbonyl ; and

alkyl is a straight or branched saturated hydrocarbon radical having from 1 to 6 carbon atoms or a cyclic saturated hydrocarbon radicals having from 3 to 6 carbon atoms ; optionally substituted on one or more carbon atoms with one or more radicals selected from the group consisting of phenyl, halo, cyano, oxo, hydroxy, formyl and amino radicals.

2. (Currently Amended) [[A]] The compound according to claim [[1]] 14,
wherein characterized in that

n — is 1 ;

m — is 1 ;

p — is 1 ;

Q is O ;

X is a covalent bond ;

each R¹ is Ar¹ or Ar¹-alkyl ;

q is 0 or 1 ;

R² is Ar² ;

Y is a covalent bond or a bivalent radical of formula -C(=O)- or -SO₂- ;

each Alk represents, independently from each other, a covalent bond; a bivalent straight or branched, saturated or unsaturated hydrocarbon radical having from 1 to 6 carbon atoms; or a cyclic saturated or unsaturated hydrocarbon radical having from 3 to 6 carbon atoms ; each radical optionally substituted on one or more carbon atoms with one or more phenyl, halo, cyano, hydroxy, formyl and amino radicals ;

L is selected from the group of hydrogen, alkyloxy, Ar³-oxy, alkyloxy-carbonyl, mono- and di(alkyl)amino, mono- and di(Ar³)amino, Ar³ and Het²;

Ar¹ is phenyl, optionally substituted with 1, 2 or 3 alkyl radicals ;

Ar² is phenyl, optionally substituted with 1, 2 or 3 alkyl radicals ;

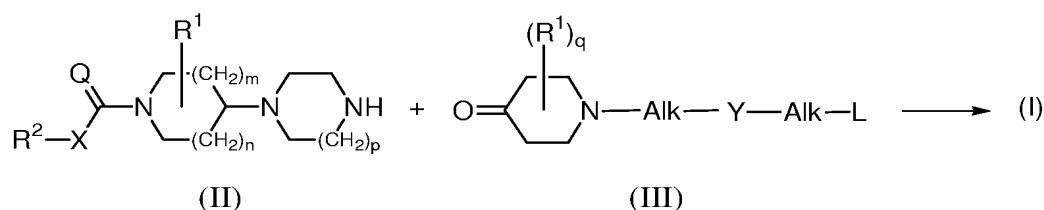
Ar³ is phenyl, optionally substituted with 1, 2 or 3 substituents each independently from each other selected from the group of alkyloxy, alkyl, halo, hydroxy, pyridinyl, morpholinyl, pyrrolidinyl, imidazo[1,2-a]pyridinyl, morpholinylcarbonyl, pyrrolidinylcarbonyl, amino and cyano ;

Het² is a monocyclic heterocyclic radical selected from the group of pyrrolidinyl, piperidinyl, morpholinyl, pyrrolyl, imidazolyl, pyrazolyl, furanyl, thienyl, isoxazolyl, thiazolyl, thiadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, and pyridazinyl ; or a bicyclic heterocyclic radical selected from the group of

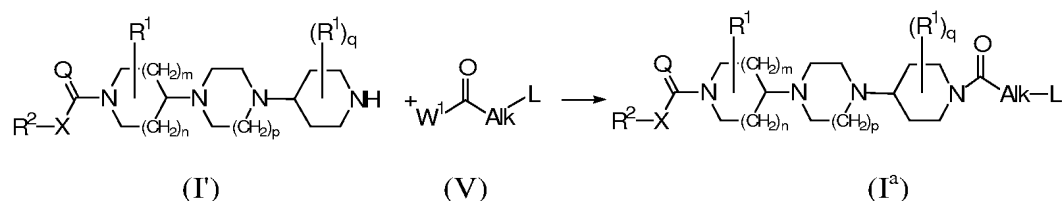
benzopiperidiny, quinolinyl, quinoxaliny, indolyl, chromenyl and benzimidazolyl ; each radical optionally substituted with one or more radicals selected from the group of Ar¹, Ar¹alkyl, halo, hydroxy, alkyl, piperidiny, pyrroly, thienyl, oxo and alkyloxycarbonyl ; and alkyl is a straight hydrocarbon radical having 1 to 6 carbon atoms, optionally substituted with one or more halo radicals .

3. (Currently Amended) [[A]] The compound according to claim [[1]]14, wherein characterized in that R¹ is Ar¹methyl and attached to the 2-position or R¹ is Ar¹ and attached to the 3-position.
4. (Currently Amended) [[A]] The compound according to claim [[1]]14, wherein characterized in that the R²-X-C(=Q)- moiety is 3,5-di-(trifluoromethyl) phenylcarbonyl.
5. (Cancelled)
6. (Currently Amended) A method of treating a warm-blooded animal suffering from a neurokinin-mediated condition comprising administering to said animal a therapeutically effective amount of a compound according to claim 1.~~The use of a compound according to claim 1 for the manufacture of a medicament for treating neurokinin mediated conditions.~~
7. (Currently Amended) The method of claim 6, wherein the neurokinin-mediated condition is use of a compound according to claim 6 for the manufacture of a medicament for~~treating~~ emesis, depression, anxiety disorder[[s]], pain, pancreatitis, micturition disorder[[s]], ~~in particular overactive bladder and~~ or irritable bowel syndrome (IBS).
8. (Previously Presented) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and, as active ingredient, a therapeutically effective amount of a compound according to claim 1.
9. (Currently Amended) [[A]] The pharmaceutical composition according to claim 8, wherein the pharmaceutical composition is an oral dosage form suitable to be orally administered.

10. (Currently Amended/Withdrawn) A process for the preparation of a composition as claimed in claim 1, wherein a pharmaceutically acceptable carrier is intimately mixed with a therapeutically effective amount of a compound as claimed in claim 1 ~~any one of claims 1-4~~.
11. (Currently Amended/Withdrawn) A process for the preparation of a compound according to ~~Formula (I)~~ claim 14, more specifically according to Formula (I^a), Formula (I^b) or Formula (I^c), wherein
- a) a final compound according to Formula (I) is obtained by reductive *N*-alkylation of an intermediate according to Formula (II) wherein R¹, R², X, Q, m, n and p are defined as in claim 14 ~~Formula (I)~~, with an *N*-substituted ~~piperidinon~~ piperidinone of Formula (III) wherein R¹, Alk, Y, L and q are defined as in claim 14 ~~Formula (I)~~, in a reaction-inert solvent and in the presence of a reducing agent ; or

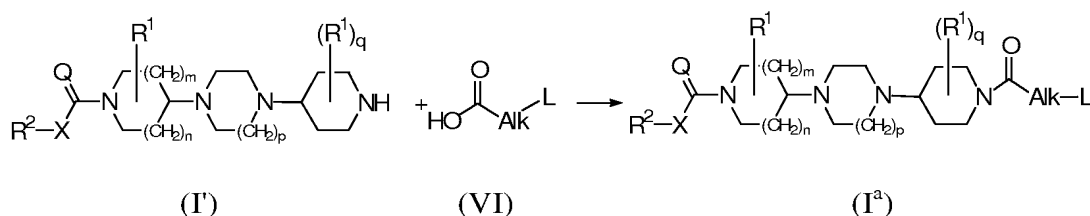


- b) a final compound according to Formula (I^a) is obtained by acylation of a final compound of Formula (I') wherein R¹, R², X, Q, m, n, p and q are defined as in claim 14 ~~Formula (I)~~, with an acyl compound of Formula (V) wherein Alk and L are defined as in claim 14 ~~Formula (I)~~ and W¹ is a leaving group, in a reaction-inert solvent and in the presence of a base ; or

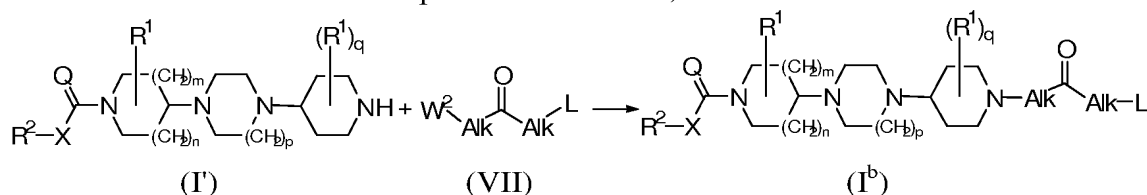


- c) a final compound according to Formula (I^a) is obtained by a base-catalyzed nucleophilic addition reaction of a final compound of Formula (I') wherein R¹, R², X, Q, m, n, p and q are defined as in claim 14 ~~Formula (I)~~, with a carboxylic acid of Formula

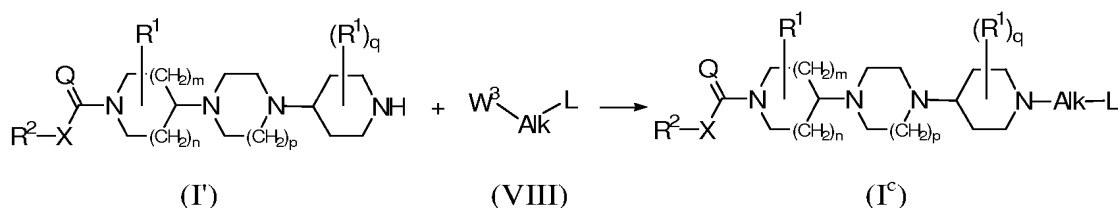
(VI) wherein Alk and L are defined as in claim 14~~Formula (I)~~, or its ester, in a reaction-inert solvent and in the presence of a base ; or



d) a final compound according to Formula (I^b) is obtained by a base-catalyzed nucleophilic addition reaction of a final compound of Formula (I') wherein R¹, R², X, Q, m, n, p and q are defined as in claim 14~~Formula (I)~~, with a compound of Formula (VIII) wherein Alk and L are defined as in claim 14~~Formula (I)~~ and W² is a leaving group, in a reaction-inert solvent and in the presence of a base ; or



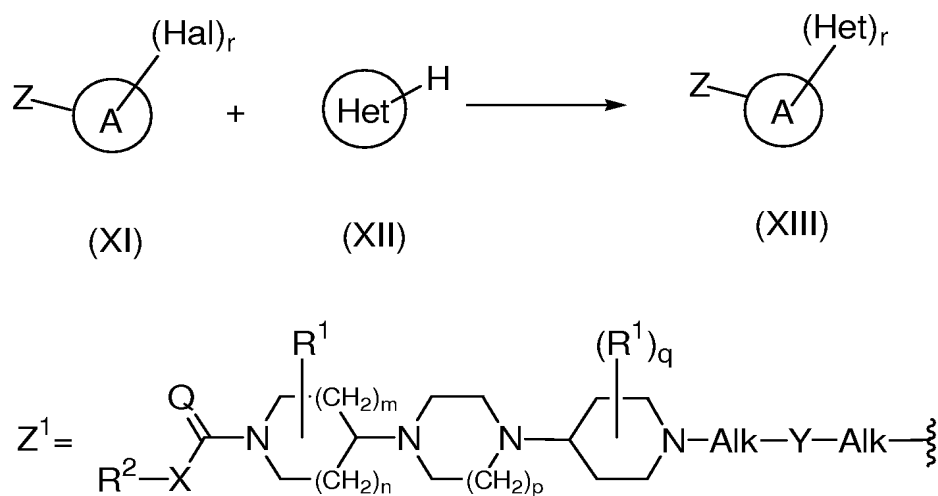
e) a final compound according to Formula (I^c) is obtained by reductive amination/alkylation of a final compound of Formula (I') wherein R¹, R², X, Q, m, n, p and q are defined as in claim 14~~Formula (I)~~ with a compound of Formula (VIII) wherein Alk and L are defined as in claim 14~~Formula (I)~~ and W³ is a leaving group, in a reaction-inert solvent and in the presence of a base ; or



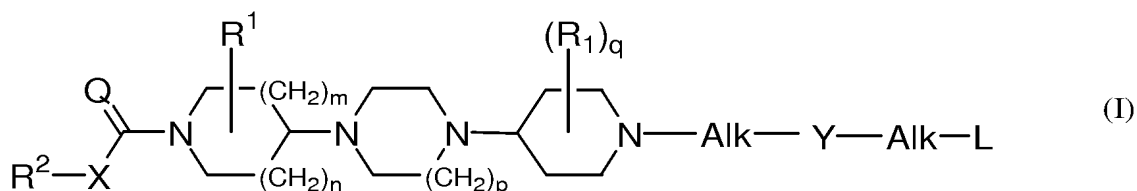
f) a final compound according to claim 14~~Formula (I)~~ is obtained by converting compounds according to claim 14 ~~Formula (I)~~ into each other following art-known transformation reactions ; and further, converting compounds according to claim 14~~Formula (I)~~ into an acid addition salt by treatment with an acid, or into a base addition

salt by treatment with a base, or conversely, the acid addition salt form may be converted into the free base by treatment with alkali, or the base addition salt may be converted into the free acid by treatment with an acid ; and by preparing the *N*-oxide and/or stereochemically isomeric forms thereof.

12. (Previously Presented/Withdrawn) A process for the preparation of a compound according to Formula (XIII), wherein a compound according to Formula (XI), wherein A is an aryl or heteroaryl, Z may be any moiety, preferably a moiety Z¹ as defined below wherein each variable is defined as in Formula (I), Hal is an halogen and r is an integer ranging from 1 to a number equal to the number of available carbon atoms in the aryl or heteroaryl-moiety A, is reacted with an unsaturated heteroaryl Het according to Formula (XII) in the presence of catalytic amounts of Pd(OAc)₂ and 1,3-bis diphenylphosphinopropane, in the presence of a suitable base, preferably Cs₂CO₃ or K(AcO) and in a reaction-inert polar solvent.



13. (Currently Amended/Withdrawn) ~~[[A]]~~The process according to claim 12, wherein Hal is bromo or iodo, A is phenyl or pyridinyl, Z is Z¹ and Het is ~~selected from the group consisting of~~imidazo[1,2-*a*]pyridinyl, pyrrolyl, or and thienyl.
14. (New) A compound according to the general Formula (I)

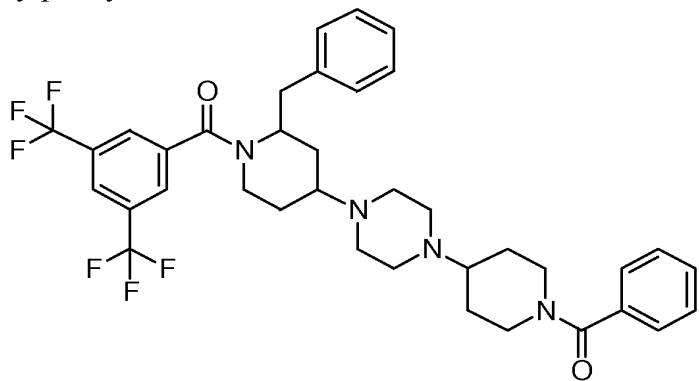


the pharmaceutically acceptable acid or base addition salts thereof and the stereochemically isomeric forms thereof, wherein:

- n is an integer equal to 1;
- m is an integer equal to 1;
- p is an integer equal to 1;
- Q is O or NR³;
- X is a covalent bond or a bivalent radical of formula -O-, -S- or -NR³-;
- each R³ independently from each other, is hydrogen or alkyl;
- each R¹ independently from each other, is selected from the group of Ar¹, Ar¹-alkyl and di(Ar¹)-alkyl;
- q is an integer equal to 0 or 1;
- R² is selected from the group consisting of alkyl, Ar², Ar²-alkyl, Het¹ or Het¹-alkyl;
- Y is a covalent bond or a bivalent radical of formula -C(=O)- or -SO₂-;
- each Alk represents, independently from each other, a covalent bond; a bivalent straight or branched, saturated or unsaturated hydrocarbon radical having from 1 to 6 carbon atoms; or a cyclic saturated or unsaturated hydrocarbon radical having from 3 to 6 carbon atoms; each radical optionally substituted on one or more carbon atoms with one or more alkyl, phenyl, halo, cyano, hydroxy, formyl and amino radicals;
- L is selected from the group of hydrogen, alkyloxy, Ar³-oxy, alkyloxycarbonyl, mono- and di(alkyl)amino, mono- and di(Ar³)amino, mono- and di(alkyloxycarbonyl)amino, Ar³, Ar³-carbonyl, Het² and Het²-carbonyl;
- Ar¹ is phenyl, optionally substituted with 1, 2 or 3 substituents each independently from each other selected from the group consisting of halo, alkyl, cyano, aminocarbonyl and alkyloxy;
- Ar² is naphthalenyl or phenyl, each optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group consisting of halo, nitro, amino, mono- and di(alkyl)amino, cyano, alkyl,

- hydroxy, alkyloxy, carboxyl, alkyloxycarbonyl, aminocarbonyl and mono- and di(alkyl)aminocarbonyl ;
- Ar³ is naphthalenyl or phenyl, optionally substituted with 1, 2 or 3 substituents each independently from each other selected from the group consisting of alkyloxy, alkyl, halo, hydroxy, pyridinyl, morpholinyl, pyrrolidinyl, imidazo[1,2-*a*]pyridinyl, morpholinylcarbonyl, pyrrolidinylcarbonyl, amino and cyano;
- Het¹ is a monocyclic heterocyclic radical selected from the group consisting of pyrrolyl, pyrazolyl, imidazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, thiadiazolyl, isothiazolyl, pyridinyl, pyrimidinyl, pyrazinyl and pyridazinyl ; or a bicyclic heterocyclic radical selected from the group of quinolinyl, quinoxalinyl, indolyl, benzimidazolyl, benzoxazolyl, benzisoxazolyl, benzothiazolyl, benzisothiazolyl, benzofuranyl and benzothienyl ; each heterocyclic radical may optionally be substituted on any atom by a radical selected from the group consisting of halo and alkyl ;
- Het² is a monocyclic heterocyclic radical selected from the group consisting of pyrrolidinyl, dioxolyl, imidazolidinyl, pyrazolidinyl, piperidinyl, morpholinyl, dithianyl, thiomorpholinyl, piperazinyl, imidazolidinyl, tetrahydrofuranyl, 2H-pyrrolyl, pyrrolinyl, imidazolinyl, pyrrazolinyl, pyrrolyl, imidazolyl, pyrazolyl, triazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, thiadiazolyl, isothiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyridazinyl and triazinyl ; or a bicyclic heterocyclic radical selected from the group consisting of benzopiperidinyl, quinolinyl, quinoxalinyl, indolyl, isoindolyl, chromenyl, benzimidazolyl, imidazo[1,2-*a*]pyridinyl, benzoxazolyl, benzisoxazolyl, benzothiazolyl, benzisothiazolyl, benzofuranyl and benzothienyl ; each radical optionally substituted with one or more radicals selected from the group consisting of Ar¹, Ar¹alkyl, halo, hydroxy, alkyl, piperidinyl, pyrrolyl, thienyl, oxo, alkyloxy, alkyloxyalkyl and alkyloxycarbonyl ; and
- alkyl is a straight or branched saturated hydrocarbon radical having from 1 to 6 carbon atoms or a cyclic saturated hydrocarbon radicals having from 3 to 6 carbon atoms ; optionally substituted on one or more carbon atoms with one or more radicals selected from the group consisting of phenyl, halo, cyano, oxo, hydroxy, formyl and amino radicals.

15. (New) The compound according to claim 14, wherein the compound of Formula I is {4-[4-(1-benzoylpiperidin-4-yl)-piperazin-1-yl]-2-benzyl-piperidin-1-yl}-(3,5-bistrifluoromethylphenyl)methanone:



16. (New) The method of claim 7, wherein the micturition disorder is overactive bladder.